

Serial No.: 10/580,202

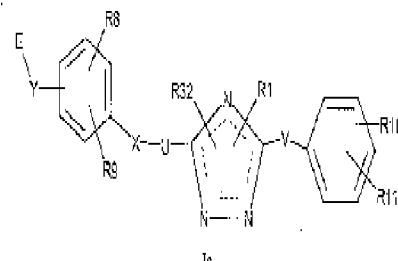
Amendments to the Claims

What is claimed is:

1. (Cancelled)

2. (Cancelled)

3. (Currently Amended) A compound wherein the compound is of the formula Ic:



and stereoisomers, or pharmaceutically acceptable salts, esters, and prodrugs thereof, wherein:

- (a) R1 is hydrogen;
- (b) R26, R27, R28 and R31 are each independently selected from the group consisting of hydrogen, hydroxy, epoxy, nitro, halo, oxo, C₁-C₃ alkyl, C₁-C₃ alkyloxy, C₁-C₃ cycloalkyl, C₁-C₃ cycloalkyloxy, C₁-C₃ cycloalkylalkyl, C₁-C₃ alkylalkyl, C(=O)R13, COOR14, OC(=O)R15, OS(=O)R16, N(R17), NR18C(=O)R19, NR18C(=O)R20, SR21, S(=O)R22, S(=O)R23, and S(=O)2R25;
- (c) R2, R3, R4, R15, R16, R17, R18, R19, R20, R21, R22, R23, R24 and R25 are each independently selected from the group consisting of hydrogen, C₁-C₃ alkyl and alkyl;
- (d) V is selected from the group consisting of C₁-C₃ alkyl;
- (e) X is selected from the group consisting of a single bond;
- (f) U is an aliphatic linker wherein one carbon atom of the aliphatic linker is substituted with from one to two substituents each independently selected from R30;

2-

=> fil cap

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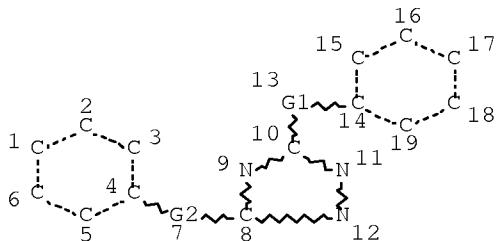
FILE COVERS 1907 - 30 Oct 2008 VOL 149 ISS 18
 FILE LAST UPDATED: 29 Oct 2008 (20081029/ED)

Caplus now includes complete International Patent Classification (IPC) reclassification data for the second quarter of 2008.

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/legal/infopolicy.html>

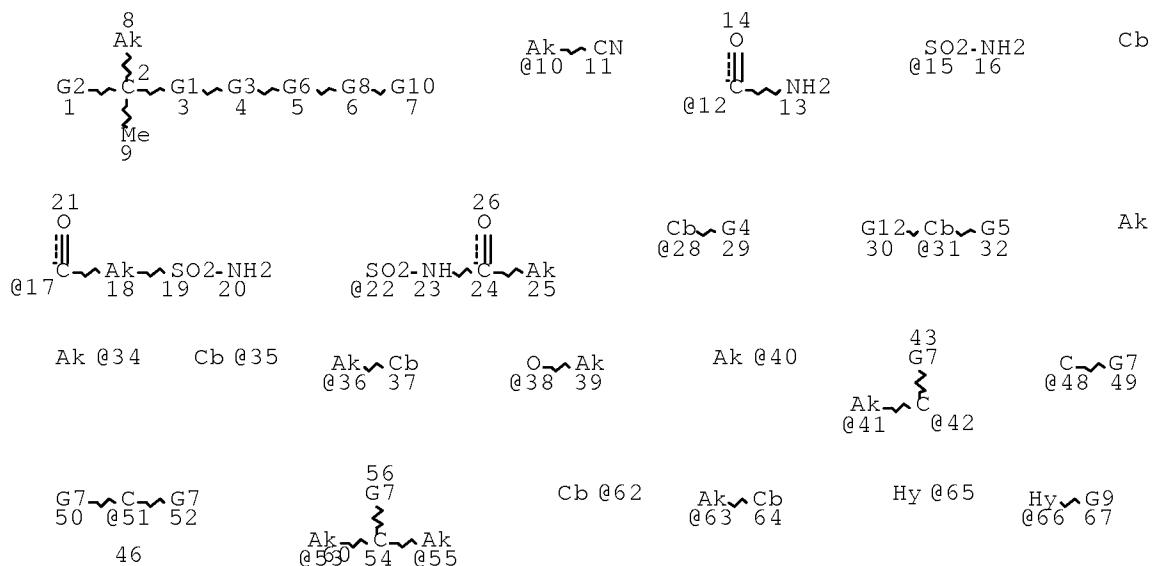
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 DEFAULT ECLEVEL IS LIMITED

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 NUMBER OF NODES IS 19

STEREO ATTRIBUTES: NONE
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 L7 STR

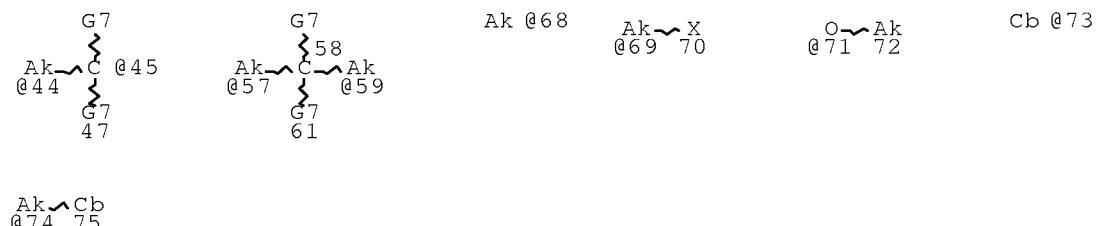


Page 1-A

@27

@33

Page 1-B



Page 2-A

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VAR G2=COOH/10/12/15/17/22

VAR G3=27/28/31

VAR G4=X/33/34/35/36/OH/38

VAR G5=34/35/36/OH/38

VAR G6=40/41-4 42-6/42-4 41-6/44-4 45-6/45-4 44-6/48/51/53-4 55-6/57-4 59-6

VAR G7=62/63

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VAR G9=X/68/69/71

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 NUMBER OF NODES IS 75

STEREO ATTRIBUTES: NONE

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FILE LAST UPDATED: 27 OCT 2008 <20081027/UP>
 MOST RECENT UPDATE: 200869 <200869/DW>
 DERWENT WORLD PATENTS INDEX SUBSCRIBER FILE, COVERS 1963 TO DATE
 >>> Now containing more than 1.2 million chemical structures in DCR <<<

>>> IPC Reform backfile reclassifications have been loaded to end of September 2008. No update date (UP) has been created for the reclassified documents, but they can be identified by 20060101/UPIC, and 20061231/UPIC, 20070601/UPIC, 20071001/UPIC, 20071130/UPIC, 20080401/UPIC, 20080701/UPIC and 20081001/UPIC.
 ECLA reclassifications to mid August and US national classification mid September 2008 have also been loaded. Update dates 20080401, 20080701 and 20081001/UPEC and /UPNC have been assigned to these. <<

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 PLEASE VISIT:

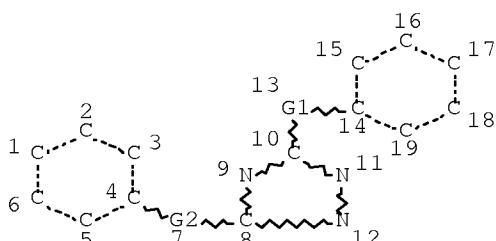
http://www.stn-international.de/training_center/patents/stn_guide.pdf

FOR DETAILS OF THE PATENTS COVERED IN CURRENT UPDATES, SEE
<http://scientific.thomsonreuters.com/support/patents/coverage/latestupdates/>

EXPLORE DERWENT WORLD PATENTS INDEX IN STN ANAVIST, VERSION 2.0:
http://www.stn-international.com/archive/presentations/DWPINAvaVist2_0608.pdf

>>> HELP for European Patent Classifications see HELP ECLA, HELP ICO <<<

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REP G1=(0-9) C

REP G2=(1-10) A

NODE ATTRIBUTES:

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DEFAULT ECLEVEL IS LIMITED

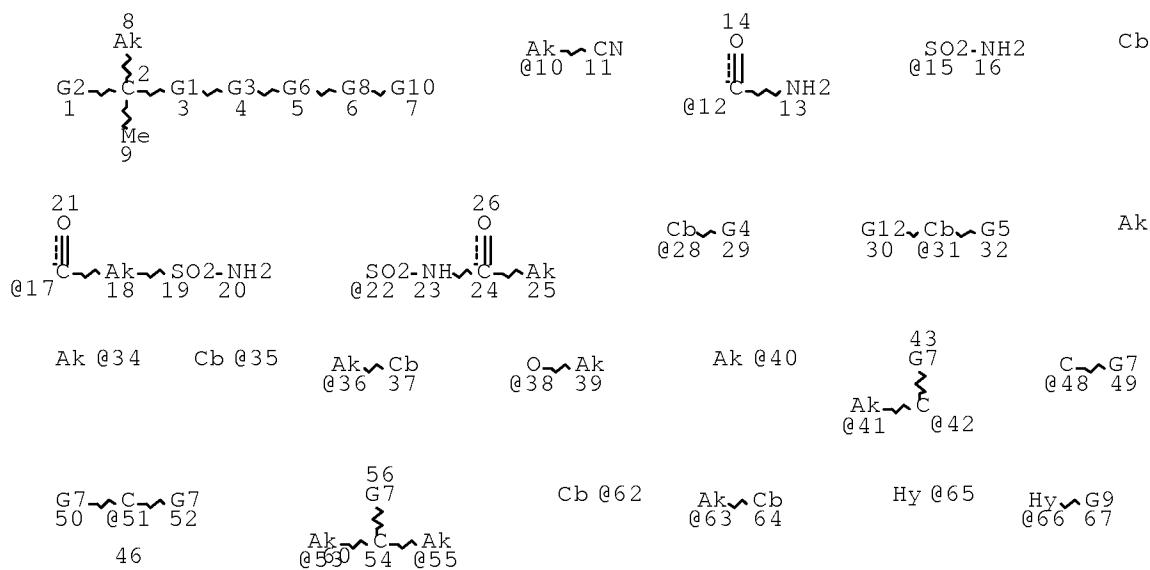
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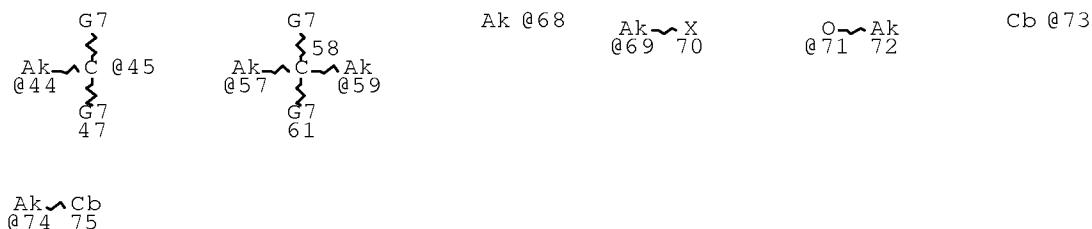


Page 1-A

@27

@33

Page 1-B



Page 2-A

VAR G1=CH2/O/S

VAR G2=COOH/10/12/15/17/22

VAR G3=27/28/31

VAR G4=X/33/34/35/36/OH/38

VAR G5=34/35/36/OH/38

VAR G6=40/41-4 42-6/42-4 41-6/44-4 45-6/45-4 44-6/48/51/53-4 55-6/57-4 59-6

VAR G7=62/63

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DEFAULT MLEVEL IS ATOM
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NUMBER OF NODES IS 75

STEREO ATTRIBUTES: NONE

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L16 1 SEA FILE=WPIX ABB=ON PLU=ON L15/DCR

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L20 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN DUPLICATE 1
ACCESSION NUMBER: 2005:638735 CAPLUS Full-text
DOCUMENT NUMBER: 143:153383
TITLE: Preparation of triazole, oxadiazole and thiadiazole derivatives as PPAR modulators for the treatment of

diabetes

INVENTOR(S): Mantlo, Nathan Bryan; Navarro, Antonio; Saeed, Ashraf; Gernert, Douglas Linn; Ma, Tianwei; Pfeifer, Lance Allen

PATENT ASSIGNEE(S): Eli Lilly and Company, USA

SOURCE: PCT Int. Appl., 175 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

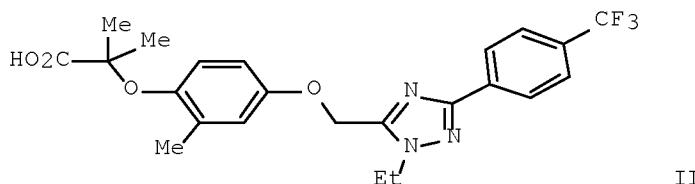
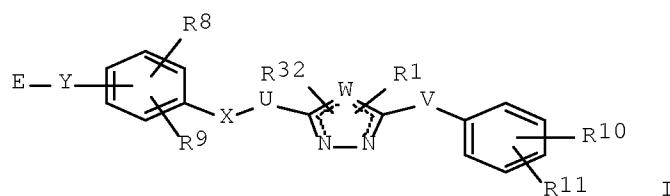
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

| PATENT NO. | KIND | DATE | APPLICATION NO. | DATE |
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| WO 2005065683 | A1 | 20050721 | WO 2004-US39775 | 20041221 |
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| RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG | | | | |
| AU 2004311909 | A1 | 20050721 | AU 2004-311909 | 20041221 |
| CA 2549385 | A1 | 20050721 | CA 2004-2549385 | 20041221 |
| EP 1725231 | A1 | 20061129 | EP 2004-812321 | 20041221 |
| R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LI, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR | | | | |
| CN 1909902 | A | 20070207 | CN 2004-80038300 | 20041221 |
| BR 2004017947 | A | 20070417 | BR 2004-17947 | 20041221 |
| JP 2007515484 | T | 20070614 | JP 2006-547018 | 20041221 |
| US 20070112045 | A1 | 20070517 | US 2006-580202 | 20060519 |
| MX 2006PA07197 | A | 20060914 | MX 2006-PA7197 | 20060622 |
| IN 2006KN01811 | A | 20070511 | IN 2006-KN1811 | 20060628 |
| PRIORITY APPLN. INFO.: | | | US 2003-532320P | P 20031222 |
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| | | | EP 2004-350159 | A 20040721 |
| | | | WO 2004-US39775 | W 20041221 |

OTHER SOURCE(S): CASREACT 143:153383; MARPAT 143:153383
GI



AB The title compds. I [X = a single bond, O, S, SO₂ and N; U = an aliphatic linker; Y = O, C, S, NH and a single bond; W = N, O or S; E = CR₃R₄A or A (wherein A = carboxy, tetrazole, alkynitrile, carboxamide, sulfonamide and acylsulfonamide; R₃ = H, alkyl, alkoxy; R₄ = H, alkyl, alkoxy, etc.; or R₃ and R₄ are optionally combined to form cycloalkyl); V = (hetero)alkyl, a bond; R₁ = H, alkyl, heteroaryl, etc.; R₈ = H, alkyl, alkenyl, halo; R₉ = H, alkyl, halo, etc.; R₁₀, R₁₁ = H, OH, CN, etc.; R₃₂ = a bond, H, halo, alkyl, etc.] which are modulators of peroxisome proliferator activated receptors (PPARs) and are useful for the treatment of diabetes and other metabolic disorders, were prepared and formulated. E.g., a multi-step synthesis of II, starting from Me glycolate and benzyl bromide, was given. The binding and cotransfection efficacy values for compds. I which are especially useful for modulating a PPAR receptor, are \leq 100 nM and \geq 50%, resp.

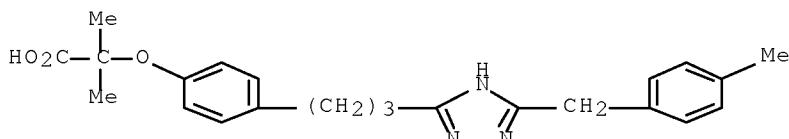
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 860262-05-1P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of triazole, oxadiazole and thiadiazole derivs. as PPAR modulators for the treatment of diabetes)

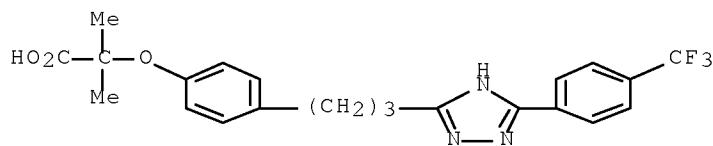
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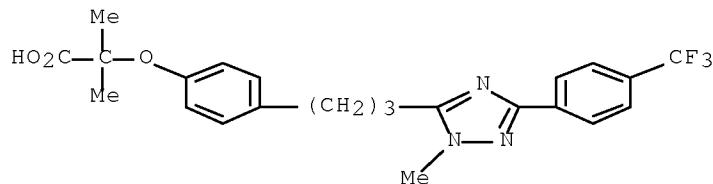
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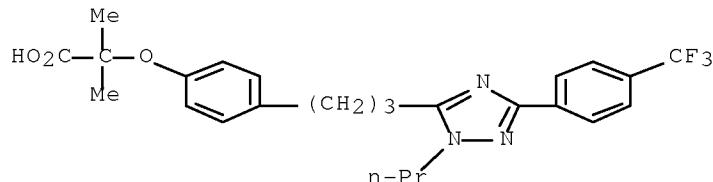
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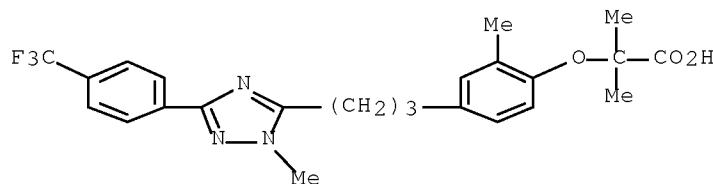
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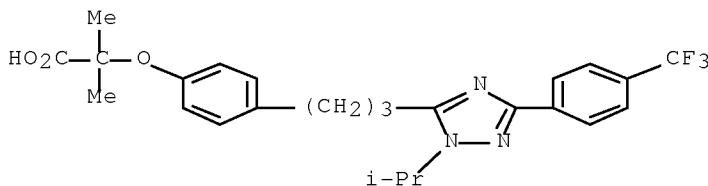
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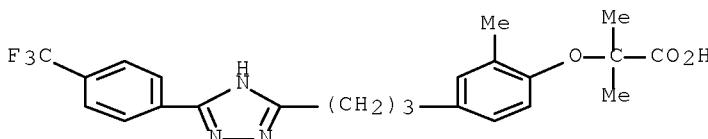
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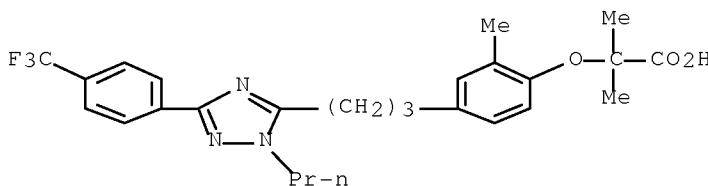
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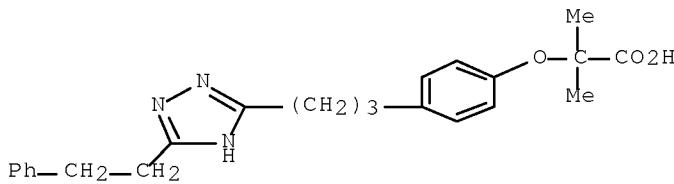
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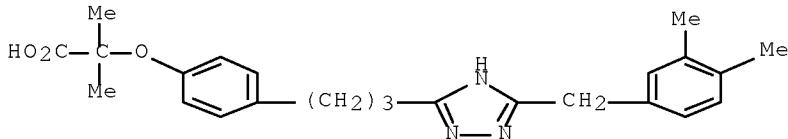


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RN 860262-05-1 CAPLUS
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 L16 1 SEA ABB=ON PLU=ON L15/DCR

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